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=> file reg COST IN U.S. DOLLARS

SINCE FILE TOTAL ENTRY SESSION 0.21 0.21

FULL ESTIMATED COST

FILE 'REGISTRY' ENTERED AT 19:33:08 ON 15 SEP 2002 USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT, PLEASE SEE "HELP USAGETERMS" FOR DETAILS. COPYRIGHT (C) 2002 American Chemical Society (ACS)

STRUCTURE FILE UPDATES: 13 SEP 2002 HIGHEST RN 450944-74-8 DICTIONARY FILE UPDATES: 13 SEP 2002 HIGHEST RN 450944-74-8

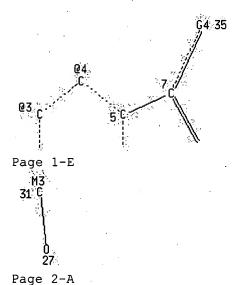
TSCA INFORMATION NOW CURRENT THROUGH MAY 20, 2002

Please note that search-term pricing does apply when conducting  ${\tt SmartSELECT}$  searches.

Crossover limits have been increased. See HELP CROSSOVER for details.

Calculated physical property data is now available. See <a href="HELP PROPERTIES">HELP PROPERTIES</a> for more information. See STNote 27, Searching Properties in the CAS Registry File, for complete details: <a href="http://www.cas.org/ONLINE/STN/STNOTES/stnotes27.pdf">http://www.cas.org/ONLINE/STN/STNOTES/stnotes27.pdf</a>

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M2 E3
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S 42 Ak 45
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Page 1-D
G1 226

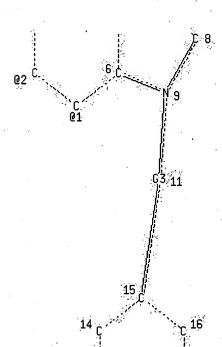






Page 2-B

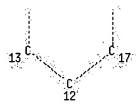
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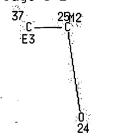
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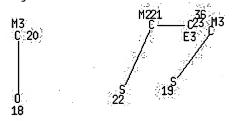
Page 3-B



Page 3-E



Page 4-D



Page 4-E

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VAR G3=42/45

VAR G4=46/47/49/50/27/28/29/30

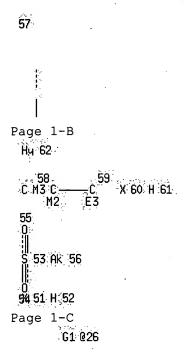
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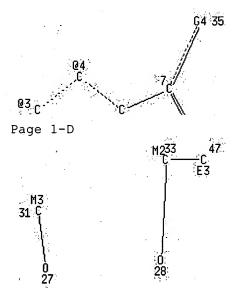
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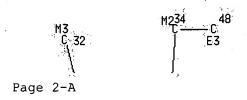
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GRAPH ATTRIBUTES:
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NUMBER OF NODES IS
STEREO ATTRIBUTES: NONE
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                                   **COMPLETE**
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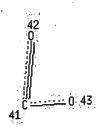
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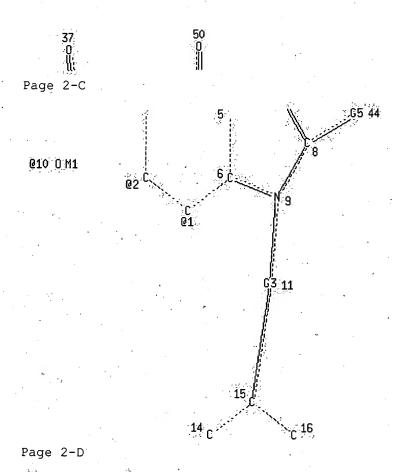
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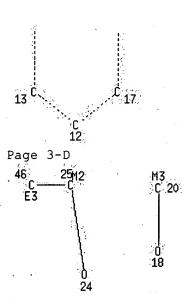






S 29 30 30 Page 3-A S 40 39 49

Page 3-C



Page 4-C M221 45 M3 C 23 M3 E3 19

VAR G1=51/52/18/19/22/24

Page 4-D

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GRAPH ATTRIBUTES:
RSPEC I
NUMBER OF NODES IS
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STEREO ATTRIBUTES: NONE
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0 ANSWERS

FULL FILE PROJECTIONS: ONLINE \*\*COMPLETE\*\*

BATCH \*\*COMPLETE\*\*

PROJECTED ITERATIONS: 318666 TO 333934

PROJECTED ANSWERS: 0 TO 0

L4 'O SEA SSS SAM L3

=> s 13 full

THE ESTIMATED SEARCH COST FOR FILE 'REGISTRY' IS 139.90 U.S. DOLLARS DO YOU WANT TO CONTINUE WITH THIS REQUEST? (Y)/N or END:y
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FULL SCREEN SEARCH COMPLETED - 323719 TO ITERATE

88.2% PROCESSED 285461 ITERATIONS

7 ANSWERS

100.0% PROCESSED 323719 ITERATIONS

7 ANSWERS

SEARCH TIME: 00.00.29

L5

7 SEA SSS FUL L3

=> file hcaplus

COST IN U.S. DOLLARS

SINCE FILE

TOTAL

ENTRY

SESSION

FULL ESTIMATED COST

149.40 149.61

FILE 'HCAPLUS' ENTERED AT 19:48:14 ON 15 SEP 2002 USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT. PLEASE SEE "HELP USAGETERMS" FOR DETAILS. COPYRIGHT (C) 2002 AMERICAN CHEMICAL SOCIETY (ACS)

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FILE COVERS 1907 - 15 Sep 2002 VOL 137 ISS 12 FILE LAST UPDATED: 13 Sep 2002 (20020913/ED)

This file contains CAS Registry Numbers for easy and accurate substance identification.

CAS roles have been modified effective December 16, 2001. Please check your SDI profiles to see if they need to be revised. For information on CAS roles, enter <a href="HELP ROLES">HELP ROLES</a> at an arrow prompt or use the CAS Roles thesaurus (/RL field) in this file.

=> s 15

1.6

4 L5

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32 KETTLE, J?/AU

L7 0 L6 AND KETTLE, J?/AU

=> s 16 and faull, a?/au

32 FAULL, A?/AU

L8 0

O L6 AND FAULL, A?/AU

=> d 16, ibib abs fhitstr, 1-4

L6 ANSWER 1 OF 4 HCAPLUS COPYRIGHT 2002 ACS



ACCESSION NUMBER:

1999:218959 HCAPLUS

DOCUMENT NUMBER:

130:325062

TITLE:

Cycloaddition reactions of furo[2,3-b]pyrroles

AUTHOR(S): Sleziak, Robert; Krutosikova, Alzbeta

CORPORATE SOURCE:

Department of Organic Chemistry, Faculty of Chemical

Technology, Slovak University of Technology,

Bratislava, SK-812 37, Slovakia

SOURCE:

Collection of Czechoslovak Chemical Communications

(1999), 64(2), 321-328

CODEN: CCCCAK; ISSN: 0010-0765

PUBLISHER: Institute of Organic Chemistry and Biochemistry,

Academy of Sciences of the Czech Republic

Journal

DOCUMENT TYPE: LANGUAGE:

English

OTHER SOURCE(S):

CASREACT 130:325062

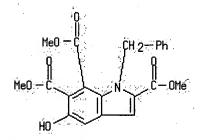
ΑB Reactions of furo[2,3-b]pyrroles (I; R = H, Me, CH2Ph, CH2OMe) with di-Me butynedioate and Et propynoate were investigated. The reaction course is influenced by the substituents on the fused system. Products of [4+2]cycloaddn. to the furan ring, leading to indole derivs., have been In the case of the reaction of I (R = H) with di-Me butynedioate, products of [4+2]cycloaddn. to the furan ring as well as of Michael addn. to the pyrrole ring, leading to N-substituted indole deriv. II have been obsd.

## IT 223715-21-7P

RL: SPN (Synthetic preparation); PREP (Preparation) (cycloaddn. reactions of furo[2,3-b]pyrroles)

HCAPLUS RN 223715-21-7

CN 1H-Indole-2,6,7-tricarboxylic acid, 5-hydroxy-1-(phenylmethyl)-, trimethyl ester (9CI) (CA INDEX NAME)



ANSWER 2 OF 4

HCAPLUS COPYRIGHT 2002 ACS



ACCESSION NUMBER:

1998:568589 HCAPLUS

DOCUMENT NUMBER:

129:175653

TITLE:

Preparation of benzenesulfonamides as elastase

inhibitors

INVENTOR(S):

Nakae, Takahiko; Kato, Masashi; Fujita, Takehito;

Kawabata, Kazuhito; Ohno, Hiroyuki

PATENT ASSIGNEE(S):

Ono Pharmaceutical Co., Ltd., Japan

SOURCE:

U.S., 150 pp.

DOCUMENT TYPE:

CODEN: USXXAM Patent

LANGUAGE:

English

FAMILY ACC. NUM. COUNT:

PATENT	INFORMATION:
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PATENT NO.	KIND	DATE	APPLICATION NO.	DATE			
US 5795890	Α	19980818	<u>US 1996-718722</u>	19960924			
JP 09165365	A2	19970624	JP 1995-272058	19950927			
JP 09278742	A2	19971028	JP 1996-271341	19960924			
JP 2881688	В2	19990412					
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AU 714025	В2	19991216					
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			JP 1996-45663 A	19960224			
		•	JP 1996-271341 A3	19960924			
·		1	US 1996-718722 A3	19960924			
OMURD GOUDON (C) MADDIM 100 175 (52							

OTHER SOURCE(S):

MARPAT 129:175653

GT

## STRUCTURE DIAGRAM TOO LARGE FOR DISPLAY - AVAILABLE VIA OFFLINE PRINT \*

The title compds. [I; R1 = C1-8 alkyl, C1-8 alkoxy, OH, etc.; n = 0-5; D =AB carbocyclic ring; R2, R3 = H, C1-4 alkyl, C1-4 alkoxy, etc.; R2R3 = C1-4 alkylidene; CR2R3 = C3-7 cycloalkyl; R4 = C1-4 alkyl, C1-4 alkoxy; two of R4, attached to the benzene nucleus at ortho positions relative to each other, represent C3-5 alkylene; m = 0-4; R5, R6 = H, OH, C1-8 alkyl, etc.; NR5R6 = heterocyclyl] and their salts, which have an inhibitory effect on elastase and therefore are useful in the prevention and/or the treatment of emphysema, rheumatoid arthritis, atherosclerosis, adult respiratory distress syndrome (ARDS), glomerular nephritis, myocardial infarction, idiopathic ulcerative colitis, and gingivitis, were prepd. and formulated. Thus, treatment of the ester II (prepn. described) with CF3CO2H in CH2Cl2/MeOPh afforded the title compd. III.HCl which showed IC50 of 0.055 μM against human polymorphonuclear elastase.

IT 190252-50-7P

RL: BAC (Biological activity or effector, except adverse); BSU (Biological study, unclassified); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)

(prepn. of benzenésulfonamidés as elastase inhibitors)

RN 190252-50-7 HCAPLUS

CN pyrrolidinyl)phenyl]butoxy]phenyl]sulfonyl]- (9CI) (CA INDEX NAME)

ANSWER 3 OF 4 HCAPLUS COPYRIGHT 2002 ACS L6



ACCESSION NUMBER: 1997:390578 HCAPLUS

DOCUMENT NUMBER: 127:5005

TITLE: Preparation of sulfamoylphenyl alkanoates as elastase

inhibitors

INVENTOR(S): Nakae, Takahiko; Kato, Masashi; Fujita, Takehito;

> Kawabata, Kazuhito; Ohno, Hiroyuki Ono Pharmaceutical Co., Ltd., Japan

SOURCE: Eur. Pat. Appl., 270 pp.

CODEN: EPXXDW

DOCUMENT TYPE:

Patent LANGUAGE: English

FAMILY ACC. NUM. COUNT:

PATENT INFORMATION:

PATENT ASSIGNEE(S):

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	JP	09278742	₹.	Ā2	19971028		JP 1996-271341	19960924			*
	JP :	2881688	•	B2	19990412					•	
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	AU	714025	•	B2	19991216						
•	ZA	9608069		A	19970520		ZA 1996-8069	19960925			
	NO	9604045		A	19970401		NO 1996-4045	19960926			
	CA .	2186665		AA	19970328	·	CA 1996-2186665	19960927			
PRIO	RITY	APPLN.	INFO.	:		Ċ	JP 1995-272058 A	19950927			
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		* · ·		1		·	TP 1996-271341 A3	19960924			

OTHER SOURCE(S): MARPAT 127:5005

GΙ

ÁΒ R1CR2R3CO2ZSO2NR5R6 [I; R1 = (un)substituted carbocyclic or heterocyclic ring; R2,R3 = H, halo, alkyl, Ph, etc.; R2R3 = alkylidene or atoms to complete a carbocyclic ring; R5,R6 = H, OH, alkyl, etc.; NR5R6 = heterocyclyl; Z = (un)substituted 1,4-phenylene] were prepd. Thus, (S) = 4 - (tert-butoxycarbonyl = 1 = pyrrolidinylsulfonyl) - 2 - methylphenol was esterified by 2-(4-pyrrolidinophenyl) butanoic acid (prepn. each given) to give title compd. II. Data for biol. activity of I were given.

IT 190252-50-7P

RL: BAC (Biological activity or effector, except adverse); BSU (Biological study, unclassified); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)

(prepn. of sulfamoylphenyl alkanoates as elastase inhibitors)

RN 190252-50-7 HCAPLUS

CN TH-Indole-2-carboxylic acid, 5-hydroxy-1-[[3-methyl-4-[1-oxo-2-[4-(1pyrrolidinyl)phenyl]butoxy]phenyl]sulfonyl]- (9CI) (CA INDEX NAME)

ANSWER 4 OF 4 HCAPLUS COPYRIGHT 2002 ACS



ACCESSION NUMBER:

DOCUMENT NUMBER:

TITLE:

AUTHOR(S):

CORPORATE SOURCE:

SOURCE:

DOCUMENT TYPE:

LANGUAGE:

GΙ

1992:651251 **HCAPLUS** 

117:251251

Substituted 4-benzylfuro[3,2-b]pyrroles

Krutosikova, Alzbeta; Hanes, Mikulas

Dep. Org. Chem., Slovak Tech. Univ., Bratislava, 812

37, Czech.

Collect. Czech. Chem. Commun. (1992), 57(7), 1487-94

CODEN: CCCCAK; ISSN: 0010-0765

Journal English

AΒ Prepn. of 4-benzylfuro[3,2-b]pyrroles is described and their reactions with selected dienophiles are discussed. Utilization of 4-acetylfuro[3,2-b]pyrroles for prepn. of 4-substituted derivs. of furo[3,2-b]pyrrole and the synthesis of Et 4-(2- and 4nitrobenzyl)furo[3,2-b]pyrrole-5-carboxylates for fusing to a 1,4-diazepine system is presented. Thus, Et furo[3,2-b]pyrrole-5carboxylate reacted with PhCH2Cl to give the benzyl deriv. I (R = CO2Et). I (R = CO2Et) was hydrolyzed to give the acid; and the acid was further decarboxylated. I (R = CO2Et, H) underwent cycloaddn. to MeO2CC≡CCO2Me to give indoledicarboxylates II.

IT 144658-74-2P

RL: SPN (Synthetic preparation); PREP (Preparation) (prepn. of)

144658-74-2 HCAPLUS RN

CN 1H-Indole-2, 4,5-tricarboxylic acid, 6-hydroxy-1-(phenylmethyl)-, 2-ethyl 4,5-dimethyl ester (9CI) (CA INDEX NAME)

=> file caold COST IN U.S. DOLLARS

SINCE FILE TOTAL ENTRY SESSION

FULL ESTIMATED COST

21.83 171.44

DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)

SINCE FILE TOTAL ENTRY SESSION

CA SUBSCRIBER PRICE

-2.48-2.48

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FILE COVERS 1907-1966 FILE LAST UPDATED: 01 May 1997 (19970501/UP)

This file contains CAS Registry Numbers for easy and accurate substance identification. Title keywords, authors, patent assignees, and patent information, e.g., patent numbers, are now searchable from 1907-1966. TIFF images of CA abstracts printed between 1907-1966 are available in the PAGE display formats.

This file supports REG1stRY for direct browsing and searching of all substance data from the REGISTRY file. Enter HELP FIRST for more information.

## => d his

L1

(FILE 'HOME' ENTERED AT 19:32:57 ON 15 SEP 2002)

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STRUCTURE UPLOADED

L2 0 S L1

L3 STRUCTURE UPLOADED

.L4 0 S L3

L5 7 S L3 FULL

FILE 'HCAPLUS' ENTERED AT 19:48:14 ON 15 SEP 2002

L6 4 S L5

L7 O S L6 AND KETTLE, J?/AU L8

0 S L6 AND FAULL, A?/AU

FILE 'CAOLD' ENTERED AT 19:49:18 ON 15 SEP 2002

s 15

L9 0 L5

=> log y

COST IN U.S. DOLLARS SINCE FILE TOTAL ENTRY SESSION FULL ESTIMATED COST 171.82

DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS) . SINCE FILE TOTAL ENTRY SESSION CA SUBSCRIBER PRICE 0.00 -2.48

STN INTERNATIONAL LOGOFF AT 19:49:31 ON 15 SEP 2002

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C:\stnweb\1.str
                      CH<sub>3</sub>
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chain nodes :
             19
    10 12
                   20
                        21
                            22
                                 23
                                       24 25 26 32 34 35 36 37 38
                                                                                  39
                                                                                       40 41 49
             56 57
                        58
                             59
         55
ring nodes :
               4 5 6 7 8
                                  9 13 14 15
                                                     16 17
     1 2
chain bonds :
     7-49 8-63 9-12 12-16 19
53-54 55-56 57-59 57-58
                          12-16 19-21 20-24 22-23 25-26 34-38 35-40 36-39 37-41 52-55
ring bonds:
     1-2 1-6 2-3 3-4 4-5 5-6 5-7 6-9 7-8 8-9 13-14 13-18 14-15 15-16 16-17
     17-18
exact/norm bonds :
     6-9 7-49 8-9 8-63 9-12 12-16 52-55 53-54 55-56 57-59 57-58
exact bonds: 5-7 7-8 19-21 20-24 22-23 25-26 34-38 35-40 36-39 37-41
normalized bonds:
_1-2 1-6 2-3 3-4 4-5 5-6 13-14 13-18 14-15 15-16 16-17 17-18
isolated ring systems :
     containing 1: 13:
G1:X,H,[*1],[*2],[*3],[*4]
G3:SO2,Ak
G4:CH3,Et,X,H,[*5],[*6],[*7],[*8]
G5:Hy,[*9],[*10]
Match level :
    1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:Atom 8:Atom 9:Atom 10:CLASS 11:CLASS 12:CLASS 13:Atom 14:Atom 15:Atom 16:Atom 17:Atom 18:Atom 19:CLASS 20:CLASS 21:CLASS 22:CLASS 23:CLASS 24:CLASS 25:CLASS 26:CLASS 32:CLASS 33 34:CLASS 35:CLASS 36:CLASS 37:CLASS 38:CLASS 39:CLASS 40:CLASS 41:CLASS 49
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34:CLASS

52:CLASS

35:CLASS

36:CLASS

26:CLASS 32:CLASS 33:CLASS 40:CLASS 41:CLASS 49:CLASS

39:CLASS 40:CLASS